### STRUCTURE UPLOADED

=> d l1 L1 HAS NO ANSWERS L1 STR

G1 H, O, F, CF2, CF3, OH

s 15 and fluoro

2489681 FLUORO

15 FLUOROS

2489681 FLUORO

(FLUORO OR FLUOROS)

1 L5 AND FLUORO

=> d 18 iall

L8

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN

RN 139601-52-8 REGISTRY

ED Entered STN: 13 Mar 1992

CN Triptolide, 16-fluoro- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

'CN Trisoxireno [4b,5:6,7:8a,9] phenanthro [1,2-c] furan, triptolide deriv.

MF C20 H23 F O6

SR CA

LC STN Files: CA, CAPLUS, TOXCENTER

DT.CA CAplus document type: Patent

RL.P Roles from patents: BIOL (Biological study); PREP (Preparation); USES (Uses)

## Ring System Data

Elemental Analysis EA	Elemental Sequence ES	the Rings SZ	RF	Identifier RID	Count
C20-C20-C20-	-=========  OC2-OC2-OC2-  OC4-C6-C6-C6	3-3-3-5-6-6-	•	6372.1.1	1

## Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	иот	ΓE
Bioconc. Factor (BCF)	+=====================================	pH 1	-===   (1)	ACD
Bioconc. Factor (BCF)	3.20	pH 4	(1)	ACD
Bioconc. Factor (BCF)	3.20	pH 7	(1)	ACD
Bioconc. Factor (BCF)	3.20	pH 8	(1)	ACD
Bioconc. Factor (BCF)	3.20	pH 10	(1)	ACD
Boiling Point (BP)	622.2+/-55.0 deg C	760.0 Torr	(1)	ACD
Enthalpy of Vap. (HVAP)	105.67+/-6.0 kJ/mol		(1)	ACD
Flash Point (FP)	330.1+/-56.7 deg C		(1)	ACD
H acceptors (HAC)	6		(1)	ACD
H donors (HD)			(1)	ACD
Koc (KOC)	80.0	pH 1	(1)	ACD
Koc (KOC)	80.0	pH 4	(1)	ACD
Koc (KOC)	80.0	pH 7	(1)	ACD
Koc (KOC)	80.0	pH 8	(1)	ACD
Koc (KOC)	80.0	pH 10	(1)	ACD
logD (LOGD)	0.97	pH 1	(1)	ACD
logD (LOGD)	0.97	pH 4	(1)	ACD
logD (LOGD)	0.97	pH 7	(1)	ACD
logD (LOGD)	0.97	pH 8	(1)	ACD
logD (LOGD)	0.97	pH 10	(1)	ACD
logP (LOGP)	0.968+/-0.754		(1)	ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1)	ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1)	ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1)	ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1)	ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1)	ACD
Molecular Weight (MW)	378.39		(1)	ACD
Vapor Pressure (VP)	4.37E-18 Torr	25.0 deg C	(1)	ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software Solaris V4.67 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

1 REFERENCES IN FILE CA (1907 TO DATE)

1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

# REFERENCE 1

ACCESSION NUMBER:

116:136234 CA

TITLE:

Isolation of 17-hydroxytriptolide and analogs as drugs

INVENTOR(S):

Ma, Pengcheng; Zheng, Jiarun; Lu, Xieyu

PATENT ASSIGNEE(S):

Chinese Academy of Medical Sciences, Institute of Skin

Disease, Peop. Rep. China

SOURCE:

Faming Zhuanli Shenqing Gongkai Shuomingshu, 12 pp.

CODEN: CNXXEV

DOCUMENT TYPE:

Patent Chinese

LANGUAGE:

INT. PATENT CLASSIF .:

MAIN:

C07D493-22

CLASSIFICATION:

63-4 (Pharmaceuticals)

Section cross-reference(s): 30

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

	_			
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
CN 1052859	A	19910710	CN 1989-105432	19891222
US 5430054	A	19950704	US 1990-629411	19901218
PRIORITY APPLN.	INFO.:		CN 1989-105432	19891222
			CN 1989-105433	19891222
	•		CN 1989-105434	19891222
•			CN 1990-105750	19901013

#### GRAPHIC IMAGE:

### ABSTRACT:

The title compds. (I; R1 = halo, OH, MeO; R2 = F, Cl, OH, R1R2 = O; R3 = halohalomethyl, CH2OH, CH2OMe, CHO, etc.), useful as antiinflammatory, antitumor, contraceptive agents, and immunosuppressants (no data), are isolated from Tripterygium wilfordii. Extraction of 20 kg T. wilfordii with 75-95% EtOH, concentration,

partition in CHCl3, and silica gel column chromatog. gave pure triptolide (I: R1R2 = 0; R3 = CH2OH), which was hydrolyzed with HX ( $\bar{X}$  = halo) to give I (R1 = X, R2 = OH, R3 = CH2OH) and further reacted to give addnl. I derivs.

SUPPL. TERM:

hydroxytriptolide isolation Tripterygium; triptolide analog

isolation Tripterygium

Ι

INDEX TERM:

Tripterygium wilfordii

(hydroxytriptolide analogs isolation from)

INDEX TERM:

139713-80-7

ROLE: PROC (Process)

(isolation of, from Tripterygium wilfordii)

INDEX TERM:

139601-39-1P 139601-41-5P 139601-40-4P 139601-42-6P 139601-43-7P 139601-44-8P 139601-45-9P 139601-46-0P

139601-47-1P, Triptolid-16-oic acid 139601-48-2P

139601-49-3P 139601-50-6P 139601-51-7P 139601-52-8P

139601-53-9P 139601-54-0P

ROLE: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); THU (Therapeutic use);